Please substitute the following amended claims for corresponding claims previously presented. A copy of the amended claims showing the requested revisions is attached.

1 (Amended). A compound of formula I,

R1 R2

R41 R41

R7 R41

R7 R44

R8 R44 R42

N N R4

R9 R44

R9 R44

 R^1 and R^2 independently represent H, $C_{1.4}$ alkyl, OR^{2b} or $N(R^{2c})R^{2d}$, or together form $O-(CH_2)_2-O-$, $-(CH_2)_3-$, $-(CH_2)_4-$ or $-(CH_2)_5-$;

R^{2b}, R^{2c} and R^{2d} independently represent H or C_{1.6} alkyl;

 R^3 represents H, $C_{1.6}$ alkyl or, together with R^4 , represents $C_{3.6}$ alkylene (which alkylene group is optionally interrupted by an O atom and/or is optionally substituted by one or more $C_{1.3}$ alkyl groups);

 R^4 represents H, $C_{1\cdot 12}$ alkyl, $C_{1\cdot 6}$ alkoxy (which latter two groups are both optionally substituted and/or terminated by one or more substituents selected

from -OH, halo, cyano, nitro, $C_{1.4}$ alkyl and/or $C_{1.4}$ alkoxy), - $(CH_2)_q$ -aryl, - $(CH_2)_q$ -oxyaryl, - $(CH_2)_q$ -Het¹ (which latter three groups are optionally substituted (at the - $(CH_2)_q$ - part and/or the aryl/Het¹ part) by one or more substituents selected from -OH, halo, cyano, nitro, - $C(O)R^{10}$, - $C(O)OR^{11}$, - $N(H)S(O)_2R^{11a}$, $C_{1.6}$ alkyl and/or $C_{1.6}$ alkoxy), - $(CH_2)_qN(H)C(O)R^8$, - $(CH_2)_qS(O)_2R^8$, - $(CH_2)_qC(O)R^8$, - $(CH_2)_qC(O)OR^8$, - $(CH_2)_qC(O)N(R^9)R^8$ or, together with R^3 , represents $C_{3.6}$ alkylene (which alkylene group is optionally interrupted by an O atom and/or is optionally substituted by one or more $C_{1.3}$ alkyl groups);

q represents 0, 1, 2, 3, 4, 5 or 6;

 R^8 represents H, $C_{1.6}$ alkyl, aryl (which latter group is optionally substituted and/or terminated by one or more substituents selected from -OH, halo, cyano, nitro, -C(O)R¹⁰, -C(O)OR¹¹, -N(H)S(O)₂R^{11a}, $C_{1.6}$ alkyl and/or $C_{1.6}$ alkoxy) or, together with R^9 , represents $C_{3.7}$ alkylene;

R⁹ represents H, C,₄ alkyl or, together with R⁸, represents C₃₋₇ alkylene;

Het¹ represents a five to twelve-membered heterocyclic ring containing one or more heteroatoms selected from oxygen, nitrogen and/or sulfur, and which also optionally includes one or more =O substituents;

R⁴¹, R⁴², R⁴³, R⁴⁴, R⁴⁵ or R⁴⁶ independently represent H or C_{1.3} alkyl;

 R^5 represents H, halo, $C_{1\cdot3}$ alkyl, $\cdot OR^{12}$, $\cdot N(R^{13})R^{12}$ or, together with R^6 , represents =0;

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 R^6 represents H, $C_{1.4}$ alkyl or, together with R^5 , represents =0;

 R^{12} represents H, $C_{1.6}$ alkyl, $\cdot S(O)_2 \cdot C_{1.4} \cdot alkyl$, $\cdot C(O)R^{14}$, $\cdot C(O)OR^{14}$, $\cdot C(O)N(R^{15})R^{15a}$ or aryl (which latter group is optionally substituted and/or terminated by one or more substituents selected from $\cdot OH$, halo, cyano, nitro, $\cdot C(O)R^{10}$, $\cdot C(O)OR^{11}$, $\cdot N(H)S(O)_2R^{11a}$, $C_{1.6}$ alkyl and/or $C_{1.6}$ alkoxy);

R¹³ represents\H or C_{1.4} alkyl;

R¹⁴ represents H or C_{1.6} alkyl;

 R^{15} and R^{15a} independently represent H or $C_{1.4}$ alkyl, or together represent $C_{3.6}$ alkylene, optionally interrupted by an O atom;

A represents a single bond, $C_{1.6}$ alkylene, $\cdot N(R^{16})(CH_2)_{r^-}$ or $\cdot O(CH_2)_{r^-}$ (in which two latter groups, the $\cdot (CH_2)_{r^-}$ group is attached to the bispidine nitrogen atom);

B represents a single bond, $C_{1\cdot4}$ alkylene, $\cdot(CH_2)_nN(R^{17})\cdot$, $\cdot(CH_2)_nS(O)_p\cdot$, $\cdot(CH_2)_nO\cdot$ (in which three latter groups, the $\cdot(CH_2)_n\cdot$ group is attached to the carbon atom bearing R^5 and R^6), $\cdot C(O)N(R^{17})\cdot$ (in which latter group, the $\cdot C(O)\cdot$ group is attached to the carbon atom bearing R^5 and R^6), $\cdot N(R^{17})C(O)O(CH_2)_n\cdot$, $\cdot N(R^{17})(CH_2)_n\cdot$ (in which two latter groups, the $N(R^{17})$ group is attached to the carbon atom bearing R^5 and R^6) or $\cdot(CH_2)_mC(H)(OH)(CH_2)_n\cdot$ (in which latter group, the $\cdot(CH_2)_m\cdot$ group is attached to the carbon atom bearing R^5 and R^6);

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m represents 1, 2 or 3;

n and r independently represent 0, 1, 2, 3 or 4;

p represents 0, 1 or 2;

R¹⁶ and R¹⁷ independently represent H or C_{1.4} alkyl;

 R^7 represents C_{16} alkyl, aryl or Het^2 , all of which groups are optionally substituted and/or terminated (as appropriate) by one or more substituents selected from -OH, cyano, halo, amino, nitro, Het^3 , -C(O) R^{10} , C(O) QR^{11} , C_{1.6} alkyl, $C_{1.6}$ alkoxy, -N(H)S(O) $_2R^{18}$, -S(O) $_2R^{19}$, -OS(O) $_2R^{20}$, -N(H)C(O)N(H) QR^{21} , -C(O)N(H) QR^{22} and/or aryl (which latter group is optionally substituted by one or more cyano groups);

Het² and Het³ independently represent a five to twelve-membered heterocyclic group containing one or more heteroatoms selected from oxygen, nitrogen and/or sulfur, and which also optionally includes one or more =0 substituents;

R¹⁸, R¹⁹ and R²⁰ independently represent C_{1.6} alkyl;

 R^{21} and R^{22} independently represent H or $C_{1.6}$ alkyl (optionally terminated by cyano); and

 R^{10} and R^{11} independently represent, at each individual occurrence, H or $C_{1.6}$ alkyl;

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 R^{1a} represents, at each individual occurrence, C_{1-6} alkyl;

or a [pharmaceutically acceptable] salt, solvate or protected derivative thereof;

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provided that:

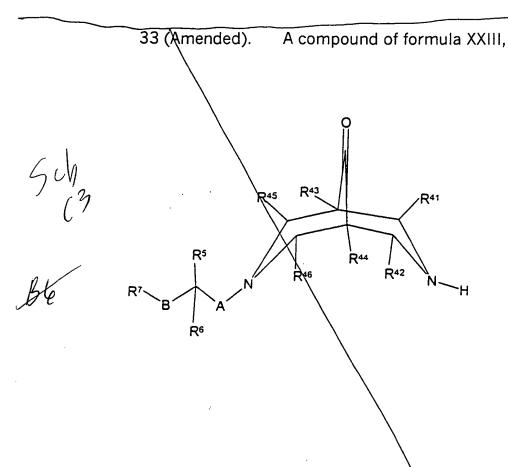
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- (a) when A and B are both single bonds and R⁷ is optionally substituted 10 aryl, then R⁵ and R⁶ do not both represent H;
- (b) when A represents a single bond, then R⁵ and R⁶ do not together represent =0; and
 - (c) when R⁵ represents -OR¹² or -N(R¹³)R¹², then:-
 - (i) A does not represent $-N(R^{16})(CH_2)_{r}$ or $O(CH_2)_{r}$; and/or
- (ii) n does not represent 0 when B represents $(CH_2)_nN(R^{17})_-$, $-(CH_2)_nN(O)_p$ or $-(CH_2)_nO_-$.

16 (Amended). A compound as claimed in Claim 15, wherein \mathbb{R}^7 represents phenyl (substituted by a cyano group and by one or more optional $C(0)N(H)\mathbb{R}^{22}$ substituent).

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26 (Amended). A compound of formula II, as defined in Claim 25, or a protected derivative thereof, provided that R^7 does not represent optionally substituted phenyl or $C_{1.6}$ alkyl.



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wherein R⁵, R⁶, R⁴¹, R⁴², R⁴³, R", R⁴⁵, R⁴⁶, A and B are as defined in Claim 1. R⁷ represents aryl or Het², all of which groups are optionally substituted and/or terminated (as appropriate) by one or more substituents selected from -OH, cyano, halo, amino, nitro, Het^3 , $-C(0)R^{10}$, $C(0)OR^{11}$, $C_{1.6}$ alkyl, $C_{1.6}$ alkoxy, $-N(H)S(O)_2R^{18}$, $-S(O)_2R^{19}$, $-OS(O)_2R^{20}$, $-N(H)C(O)N(H)R^{21}$, $-C(O)N(H)R^{22}$ and/or aryl (which latter group is optionally substituted by one or more cyano groups); or a protected derivative thereof.

Please add the following new claims.

38 (New). A compound as claimed in Claim 16, wherein the cyano group is in the 4-position relative to B.

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 β 7 39 (New). A method as claimed in Claim 24, wherein the arrhythmia is an atrial or a ventricular arrhythmia.